Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051117\
Data File : BF095107.D
Acq On : 12 May 2017 9:07
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 12 13:55:38 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 03 17:24:51 2017
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	115	-0.03
2		1,4-Dioxane	0.588	0.580	1.4	120	-0.08
3		Pyridine	1.364	1.250	8.4	109	-0.07
4		n-Nitrosodimethylamine	0.568	0.491	13.6	104	-0.06
5	S	2-Fluorophenol	1.200	1.147	4.4	111	-0.04
6		Aniline	1.817	1.942	-6.9	118	-0.02
7	S	Phenol-d6	1.439	1.470	-2.2	118	-0.02
8		2-Chlorophenol	1.365	1.428	-4.6	122	-0.02
9		Benzaldehyde	0.879	0.871	0.9	113	-0.02
10	C	Phenol	1.517	1.506	0.7	115	-0.02
11		bis(2-Chloroethyl)ether	1.252	1.269	-1.4	117	-0.03
12		1,3-Dichlorobenzene	1.549	1.553	-0.3	124	-0.04
13	C	1,4-Dichlorobenzene	1.571	1.583	-0.8	116	-0.03
14		1,2-Dichlorobenzene	1.466	1.491	-1.7	119	-0.04
15		Benzyl Alcohol	0.926	1.071	-15.7	128	-0.02
16		2,2'-oxybis(1-Chloropropane	1.772	1.694	4.4	114	-0.02
17		2-Methylphenol	1.117	1.122	-0.4	121	-0.02
18		Hexachloroethane	0.532	0.509	4.3	109	-0.04
19	P	n-Nitroso-di-n-propylamine	0.973	0.990	-1.7	120	-0.02
20		3+4-Methylphenols	1.518	1.546	-1.8	118	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	112	-0.03
22		Acetophenone	0.480	0.488	-1.7	116	-0.02
23	S	Nitrobenzene-d5	0.317	0.330	-4.1	116	-0.02
24		Nitrobenzene	0.332	0.334	-0.6	116	-0.02
25		Isophorone	0.566	0.592	-4.6	118	-0.02
26	C	2-Nitrophenol	0.179	0.182	-1.7	112	-0.02
27		2,4-Dimethylphenol	0.290	0.299	-3.1	120	-0.02
28		bis(2-Chloroethoxy)methane	0.392	0.418	-6.6	121	-0.02
29	C	2,4-Dichlorophenol	0.274	0.288	-5.1	123	-0.02
30		1,2,4-Trichlorobenzene	0.293	0.300	-2.4	118	-0.02
31		Naphthalene	1.005	1.021	-1.6	117	-0.03
32		Benzoic acid	0.177	0.038	78.5#	25#	-0.05
33		4-Chloroaniline	0.375	0.373	0.5	112	-0.02
34	C	Hexachlorobutadiene	0.155	0.143	7.7	106	-0.04
35		Caprolactam	0.082	0.086	-4.9	114	0.00
36	C	4-Chloro-3-methylphenol	0.293	0.311	-6.1	120	-0.01
37		2-Methylnaphthalene	0.660	0.692	-4.8	121	-0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	121	-0.03
39		1,2,4,5-Tetrachlorobenzene	0.615	0.599	2.6	112	-0.03
40	P	Hexachlorocyclopentadiene	0.221	0.161	27.1#	81	-0.04
41	S	2,4,6-Tribromophenol	0.164	0.153	6.7	106	-0.03
42	C	2,4,6-Trichlorophenol	0.411	0.425	-3.4	120	-0.02
43		2,4,5-Trichlorophenol	0.441	0.404	8.4	105	0.00
44	S	2-Fluorobiphenyl	1.390	1.305	6.1	111	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051117\
Data File : BF095107.D
Acq On : 12 May 2017 9:07
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 12 13:55:38 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 03 17:24:51 2017
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.684	1.640	2.6	112	-0.03
46		2-Chloronaphthalene	1.360	1.227	9.8	107	-0.02
47		2-Nitroaniline	0.372	0.362	2.7	116	-0.02
48		Acenaphthylene	2.130	2.060	3.3	114	-0.03
49		Dimethylphthalate	1.642	1.598	2.7	114	-0.03
50		2,6-Dinitrotoluene	0.335	0.342	-2.1	118	-0.02
51	C	Acenaphthene	1.272	1.248	1.9	116	-0.03
52		3-Nitroaniline	0.360	0.389	-8.1	125	0.00
53	P	2,4-Dinitrophenol	0.156	0.008#	94.9#	6#	0.06
54		Dibenzofuran	1.760	1.876	-6.6	128	-0.03
55	P	4-Nitrophenol	0.216	0.167	22.7	85	0.02
56		2,4-Dinitrotoluene	0.430	0.482	-12.1	128	0.00
57		Fluorene	1.531	1.466	4.2	117	-0.03
58		2,3,4,6-Tetrachlorophenol	0.278	0.276	0.7	118	-0.02
59		Diethylphthalate	1.574	1.588	-0.9	123	-0.03
60		4-Chlorophenyl-phenylether	0.673	0.676	-0.4	118	-0.03
61		4-Nitroaniline	0.330	0.354	-7.3	128	0.00
62		Azobenzene	1.265	1.305	-3.2	119	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	111	-0.02
64		4,6-Dinitro-2-methylphenol	0.134	0.028	79.1#	23#	0.00
65	c	n-Nitrosodiphenylamine	0.726	0.757	-4.3	117	-0.03
66		4-Bromophenyl-phenylether	0.211	0.209	0.9	108	-0.04
67		Hexachlorobenzene	0.217	0.216	0.5	111	-0.02
68		Atrazine	0.205	0.209	-2.0	119	-0.02
69	C	Pentachlorophenol	0.085	0.046	45.9#	63	-0.01
70		Phenanthrene	1.149	1.142	0.6	113	-0.03
71		Anthracene	1.174	1.203	-2.5	115	-0.02
72		Carbazole	1.068	1.161	-8.7	124	-0.02
73		Di-n-butylphthalate	1.332	1.571	-17.9	130	-0.03
74	C	Fluoranthene	1.179	1.255	-6.4	118	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	100	-0.02
76		Benzidine	0.465	0.357	23.2	67	-0.01
77		Pyrene	1.496	1.651	-10.4	109	-0.02
78	S	Terphenyl-d14	0.908	0.953	-5.0	105	-0.03
79		Butylbenzylphthalate	0.696	0.860	-23.6	121	-0.03
80		Benzo(a)anthracene	1.164	1.204	-3.4	103	-0.02
81		3,3'-Dichlorobenzidine	0.402	0.362	10.0	86	-0.02
82		Chrysene	1.125	1.161	-3.2	102	-0.02
83		Bis(2-ethylhexyl)phthalate	0.991	1.220	-23.1	119	-0.04
84	c	Di-n-octyl phthalate	1.625	1.865	-14.8	111	-0.04
85		Indeno(1,2,3-cd)pyrene	0.914	1.119	-22.4	125	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	113	-0.02
87		Benzo(b)fluoranthene	1.236	1.145	7.4	96	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051117\
Data File : BF095107.D
Acq On : 12 May 2017 9:07
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 12 13:55:38 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 03 17:24:51 2017
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.192	1.235	-3.6	114	-0.02
89 C	Benzo(a)pyrene	1.140	1.109	2.7	105	-0.02
90	Dibenzo(a,h)anthracene	0.942	1.050	-11.5	124	-0.02
91	Benzo(g,h,i)perylene	0.924	1.080	-16.9	133	-0.02

(#) = Out of Range SPCC's out = 1 CCC's out = 1